wwPDB X-ray Structure Validation Summary Report (i)

Jan 4, 2024 – 02:39 am GMT

PDB ID	:	5FSA
Title	:	Crystal structure of sterol 14-alpha demethylase (CYP51) from a pathogenic yeast Candida albicans in complex with the antifungal drug posaconazole
Authors	:	Hargrove, T.Y.; Wawrzak, Z.; Friggeri, L.; Lepesheva, G.I.
Deposited on	:	2016-01-02
Resolution	:	2.86 Å(reported)

This is a wwPDB X-ray Structure Validation Summary Report for a publicly released PDB entry.

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

The following versions of software and data (see references (i)) were used in the production of this report:

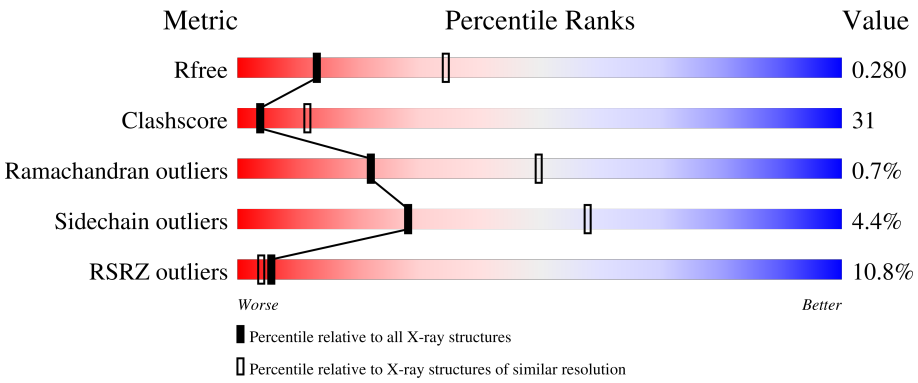
MolProbity	:	4.02b-467
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.86 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	3168 (2.90-2.82)
Clashscore	141614	3438 (2.90-2.82)
Ramachandran outliers	138981	3348 (2.90-2.82)
Sidechain outliers	138945	3351 (2.90-2.82)
RSRZ outliers	127900	3103 (2.90-2.82)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=5% The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	490	9% 60% 36% • •
1	B	490	12% 57% 37% 5% •

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 8163 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

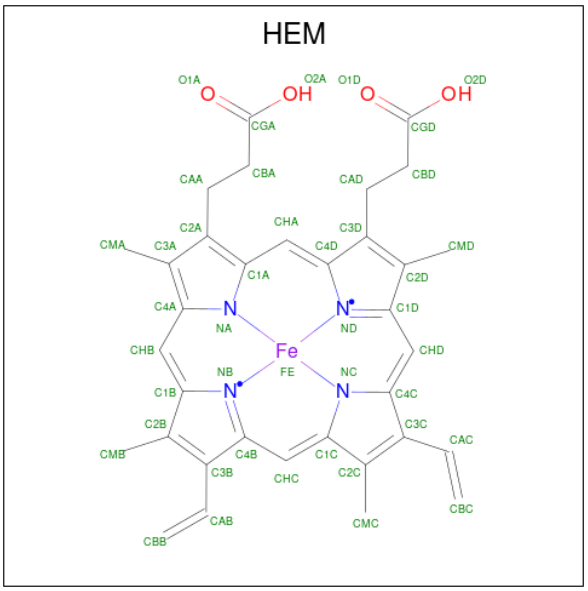
- Molecule 1 is a protein called CYP51 VARIANT1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	484	3929	2539	652	722	16	0	0	0
1	B	483	3919	2532	650	721	16	0	0	0

There are 22 discrepancies between the modelled and reference sequences:

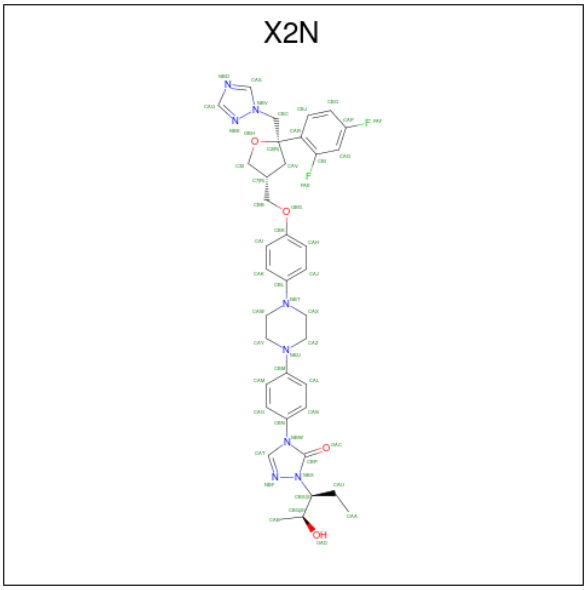
Chain	Residue	Modelled	Actual	Comment	Reference
A	43	MET	-	expression tag	UNP Q9P4W0
A	44	ALA	-	expression tag	UNP Q9P4W0
A	45	LYS	-	expression tag	UNP Q9P4W0
A	46	LYS	-	expression tag	UNP Q9P4W0
A	47	THR	-	expression tag	UNP Q9P4W0
A	48	PRO	-	expression tag	UNP Q9P4W0
A	529	HIS	-	expression tag	UNP Q9P4W0
A	530	HIS	-	expression tag	UNP Q9P4W0
A	531	HIS	-	expression tag	UNP Q9P4W0
A	532	HIS	-	expression tag	UNP Q9P4W0
A	263	LEU	SER	engineered mutation	UNP Q9P4W0
B	43	MET	-	expression tag	UNP Q9P4W0
B	44	ALA	-	expression tag	UNP Q9P4W0
B	45	LYS	-	expression tag	UNP Q9P4W0
B	46	LYS	-	expression tag	UNP Q9P4W0
B	47	THR	-	expression tag	UNP Q9P4W0
B	48	PRO	-	expression tag	UNP Q9P4W0
B	529	HIS	-	expression tag	UNP Q9P4W0
B	530	HIS	-	expression tag	UNP Q9P4W0
B	531	HIS	-	expression tag	UNP Q9P4W0
B	532	HIS	-	expression tag	UNP Q9P4W0
B	263	LEU	SER	engineered mutation	UNP Q9P4W0

- Molecule 2 is PROTOPORPHYRIN IX CONTAINING FE (three-letter code: HEM) (formula: $C_{34}H_{32}FeN_4O_4$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	Fe	N	O		
2	A	1	43	34	1	4	4	0	0
			Total	C	Fe	N	O		
2	B	1	43	34	1	4	4	0	0

- Molecule 3 is POSACONAZOLE (three-letter code: X2N) (formula: $C_{37}H_{42}F_2N_8O_4$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	F	N	O		
3	A	1	51	37	2	8	4	0	0
			Total	C	F	N	O		
3	B	1	51	37	2	8	4	0	0

- Molecule 4 is water.

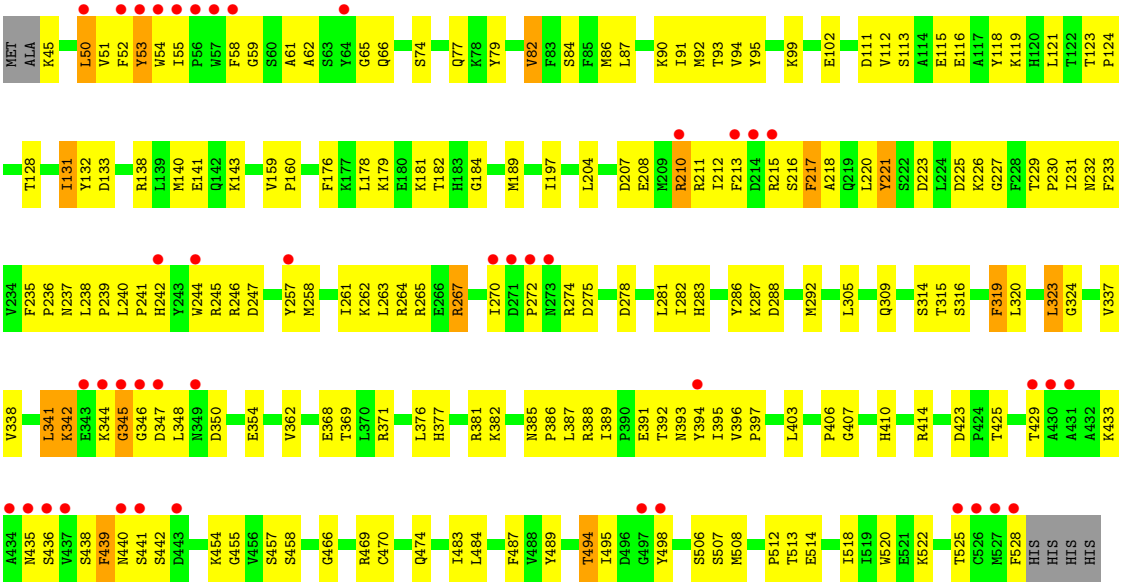
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
4	A	61	61	61	0	0
			Total	O		
4	B	66	66	66	0	0

3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density (RSRZ > 2). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

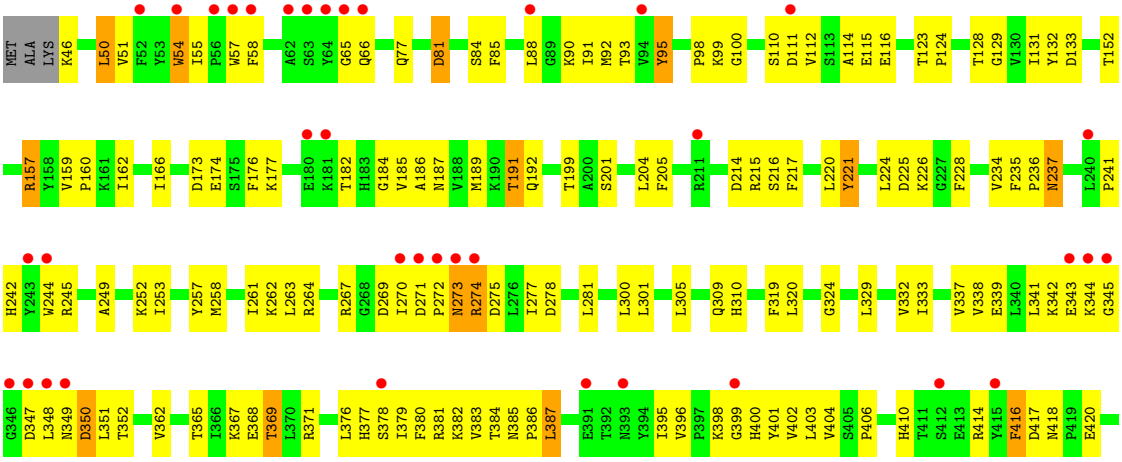
• Molecule 1: CYP51 VARIANT1

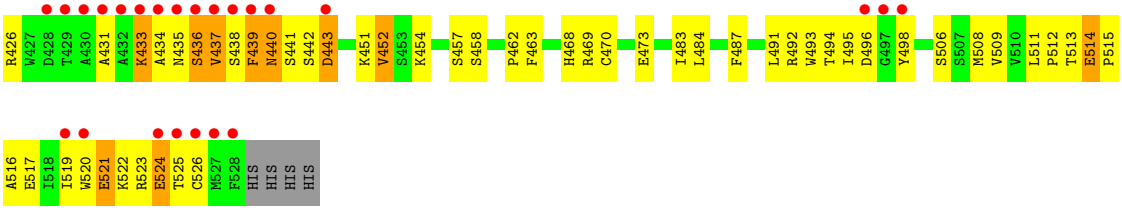
Chain A:



• Molecule 1: CYP51 VARIANT1

Chain B:




R426
W427
D428
T429
A430
A431
A432
K433
A434
N435
S436
V437
S438
F439
N440
S441
S442
D443
K451
V452
S453
K454
S457
S458
P462
F463
H468
R469
C470
E473
I483
L484
F487
L491
R492
W493
T494
I495
D496
G497
Y498
S506
S507
M508
V509
V510
L511
P512
T513
E514
P515

A516
E517
I518
I519
W520
E521
K522
R523
E524
T525
C526
M527
F528
HIS
HIS
HIS
HIS

4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	179.59Å 72.91Å 79.71Å 90.00° 96.13° 90.00°	Depositor
Resolution (Å)	50.39 – 2.86 50.39 – 2.86	Depositor EDS
% Data completeness (in resolution range)	99.3 (50.39-2.86) 99.3 (50.39-2.86)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.18 (at 2.86Å)	Xtriage
Refinement program	REFMAC 5.8.0135	Depositor
R, R_{free}	0.241 , 0.256 0.243 , 0.280	Depositor DCC
R_{free} test set	1235 reflections (5.20%)	wwPDB-VP
Wilson B-factor (Å^2)	66.1	Xtriage
Anisotropy	0.142	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.32 , 63.9	EDS
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.91	EDS
Total number of atoms	8163	wwPDB-VP
Average B, all atoms (Å^2)	73.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 11.07% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality (i)

5.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section: X2N, HEM

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\#\|Z\| >5$	RMSZ	$\#\|Z\| >5$
1	A	0.27	0/4040	0.56	1/5475 (0.0%)
1	B	0.31	0/4030	0.62	2/5462 (0.0%)
All	All	0.29	0/8070	0.59	3/10937 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand.A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	514	GLU	C-N-CD	5.73	140.44	128.40
1	A	345	GLY	N-CA-C	-5.64	99.01	113.10
1	B	343	GLU	N-CA-C	5.04	124.62	111.00

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	342	LYS	Peptide

5.2 Too-close contacts (i)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3929	0	3865	205	1
1	B	3919	0	3847	285	0
2	A	43	0	30	5	0
2	B	43	0	30	10	0
3	A	51	0	42	5	0
3	B	51	0	42	6	0
4	A	61	0	0	2	0
4	B	66	0	0	3	0
All	All	8163	0	7856	494	1

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 31.

The worst 5 of 494 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:272:PRO:HB2	1:A:274:ARG:HG3	1.32	1.09
1:B:523:ARG:HG3	1:B:524:GLU:H	1.25	1.00
1:B:506:SER:HB2	3:B:590:X2N:HAT	1.47	0.97
1:B:263:LEU:O	1:B:267:ARG:HG2	1.64	0.96
1:B:184:GLY:HA2	1:B:521:GLU:HB3	1.45	0.96

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:53:TYR:OH	1:A:237:ASN:O[2_756]	2.03	0.17

5.3 Torsion angles (i)

5.3.1 Protein backbone (i)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	482/490 (98%)	455 (94%)	25 (5%)	2 (0%)	34	62
1	B	481/490 (98%)	448 (93%)	28 (6%)	5 (1%)	15	40
All	All	963/980 (98%)	903 (94%)	53 (6%)	7 (1%)	22	50

5 of 7 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	524	GLU
1	A	342	LYS
1	B	440	ASN
1	B	216	SER
1	B	350	ASP

5.3.2 Protein sidechains (i)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	431/436 (99%)	416 (96%)	15 (4%)	36	67
1	B	430/436 (99%)	407 (95%)	23 (5%)	22	50
All	All	861/872 (99%)	823 (96%)	38 (4%)	28	58

5 of 38 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	387	LEU

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	443	ASP
1	B	414	ARG
1	B	436	SER
1	B	521	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 6 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	273	ASN
1	B	418	ASN
1	B	468	HIS
1	B	187	ASN
1	A	187	ASN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains (i)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

4 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
3	X2N	B	590	2	51,57,57	1.87	8 (15%)	57,82,82	2.33	22 (38%)
3	X2N	A	590	2	51,57,57	1.91	8 (15%)	57,82,82	2.22	19 (33%)
2	HEM	B	580	1,3	41,50,50	1.73	11 (26%)	45,82,82	2.35	16 (35%)
2	HEM	A	580	1,3	41,50,50	1.86	12 (29%)	45,82,82	1.99	14 (31%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	X2N	B	590	2	-	9/38/59/59	0/7/7/7
3	X2N	A	590	2	-	10/38/59/59	0/7/7/7
2	HEM	B	580	1,3	-	2/12/54/54	-
2	HEM	A	580	1,3	-	0/12/54/54	-

The worst 5 of 39 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	590	X2N	CBN-NBW	-6.81	1.34	1.44
3	B	590	X2N	CAR-CBI	6.70	1.48	1.38
3	A	590	X2N	CAR-CBI	6.64	1.48	1.38
3	B	590	X2N	CBN-NBW	-6.48	1.34	1.44
3	A	590	X2N	CBP-NBW	-5.38	1.33	1.39

The worst 5 of 71 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	B	580	HEM	CHC-C4B-NB	6.57	131.57	124.43
3	B	590	X2N	CAS-NBD-CAQ	6.45	109.61	102.34
3	A	590	X2N	CAS-NBD-CAQ	6.38	109.53	102.34
3	B	590	X2N	CAO-CBN-NBW	5.55	126.81	119.63
3	B	590	X2N	CBS-NBX-CBP	5.35	135.79	125.09

There are no chirality outliers.

5 of 21 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	590	X2N	CAV-C7-CBB-OBG
3	A	590	X2N	CAR-C2-CBC-NBV

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms
3	A	590	X2N	OBH-C2-CBC-NBV
3	B	590	X2N	CAV-C7-CBB-OBG
3	B	590	X2N	CAR-C2-CBC-NBV

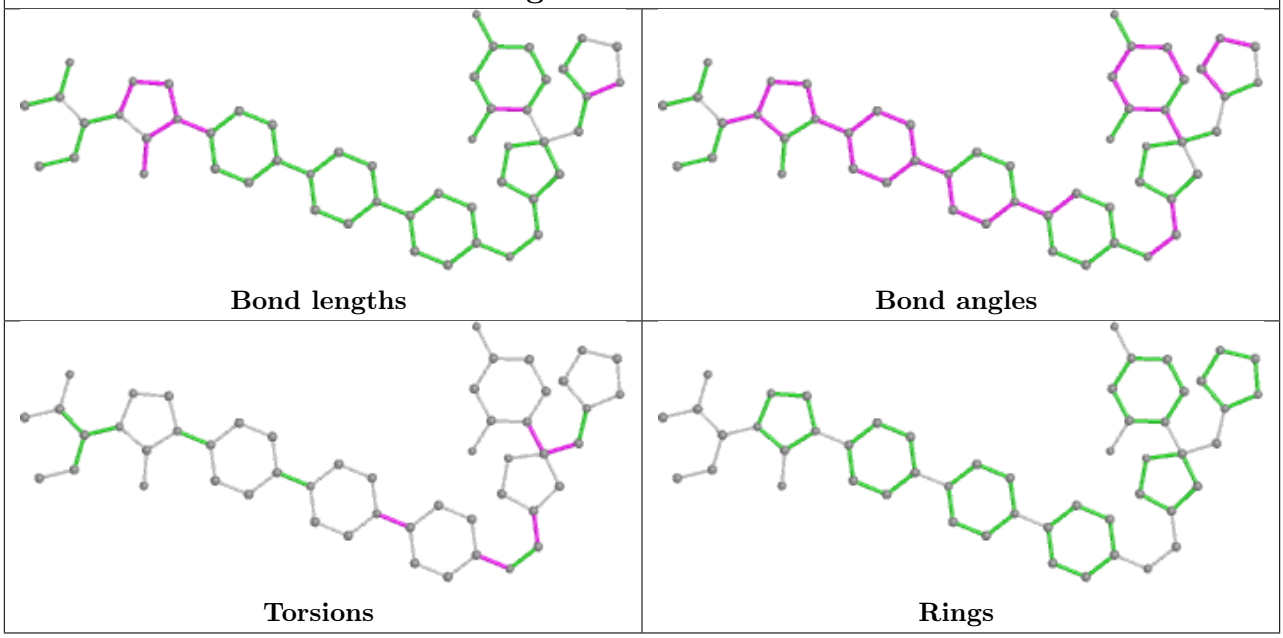
There are no ring outliers.

4 monomers are involved in 26 short contacts:

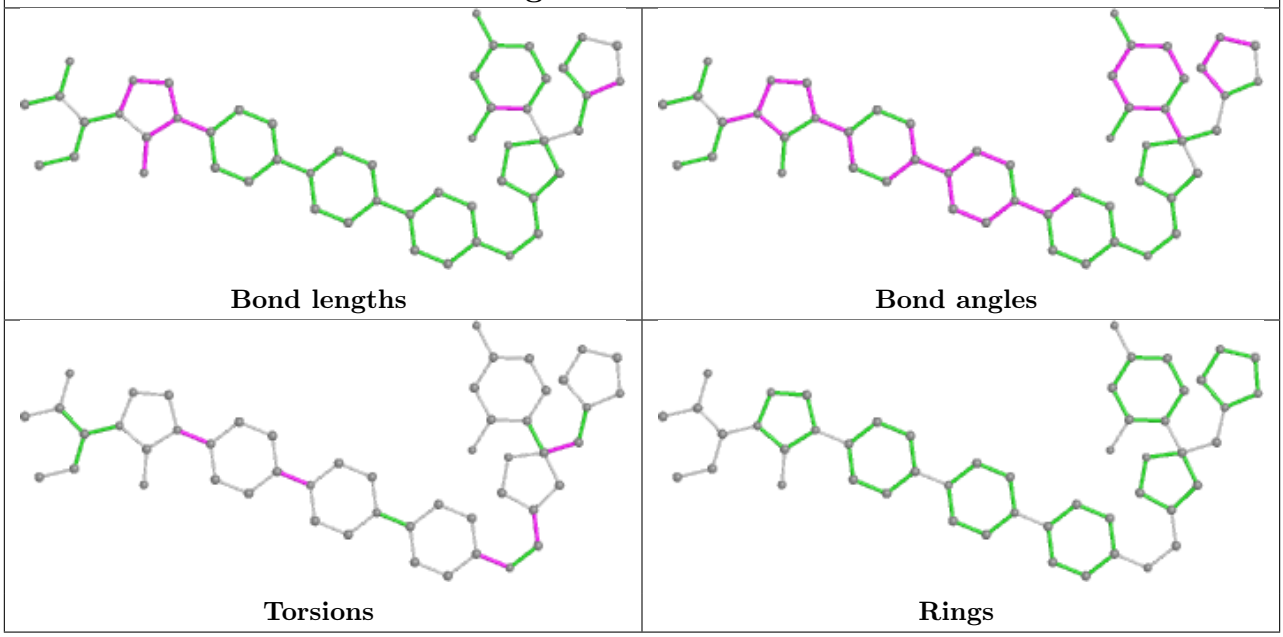
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	590	X2N	6	0
3	A	590	X2N	5	0
2	B	580	HEM	10	0
2	A	580	HEM	5	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less then 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

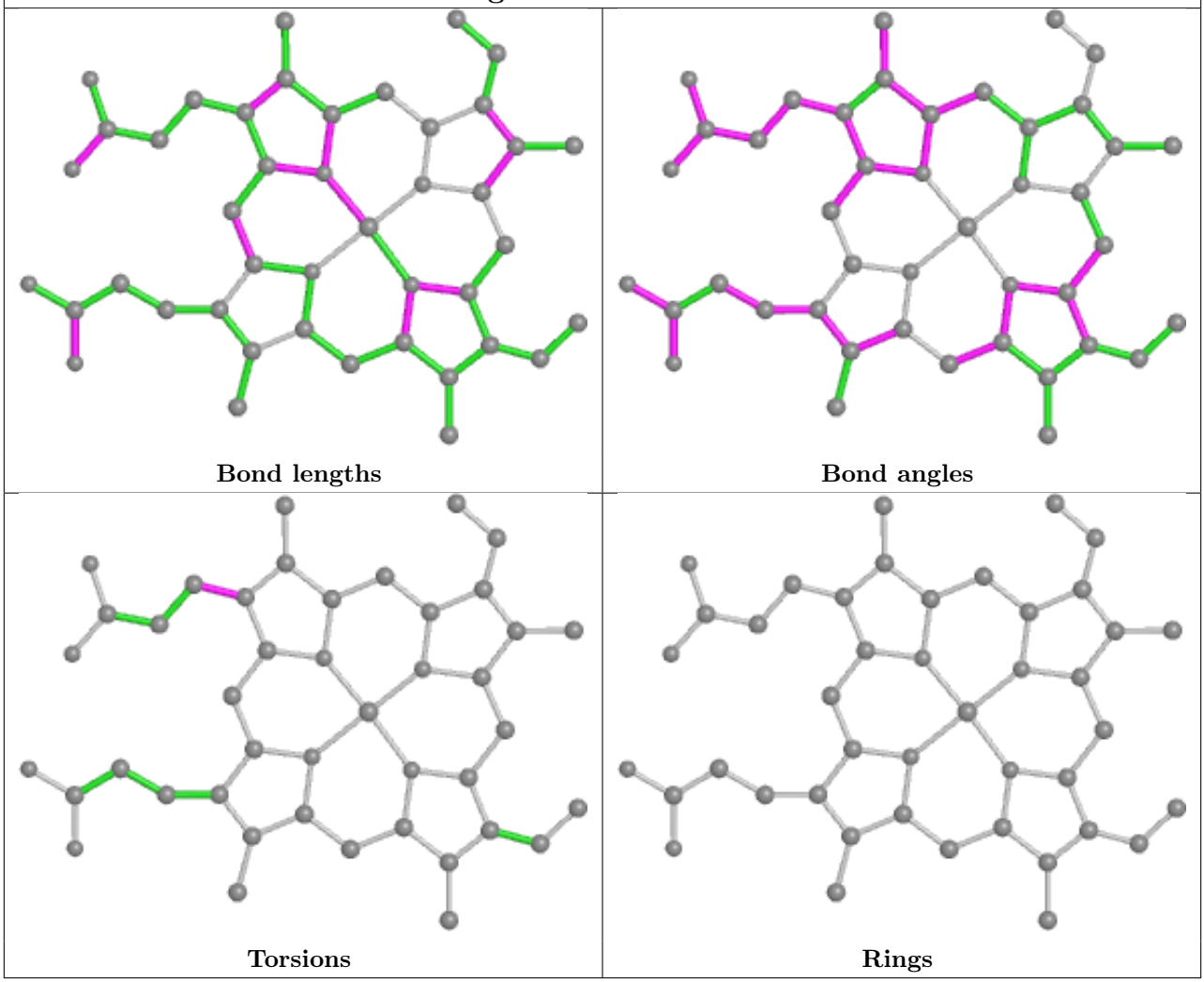
Ligand X2N B 590



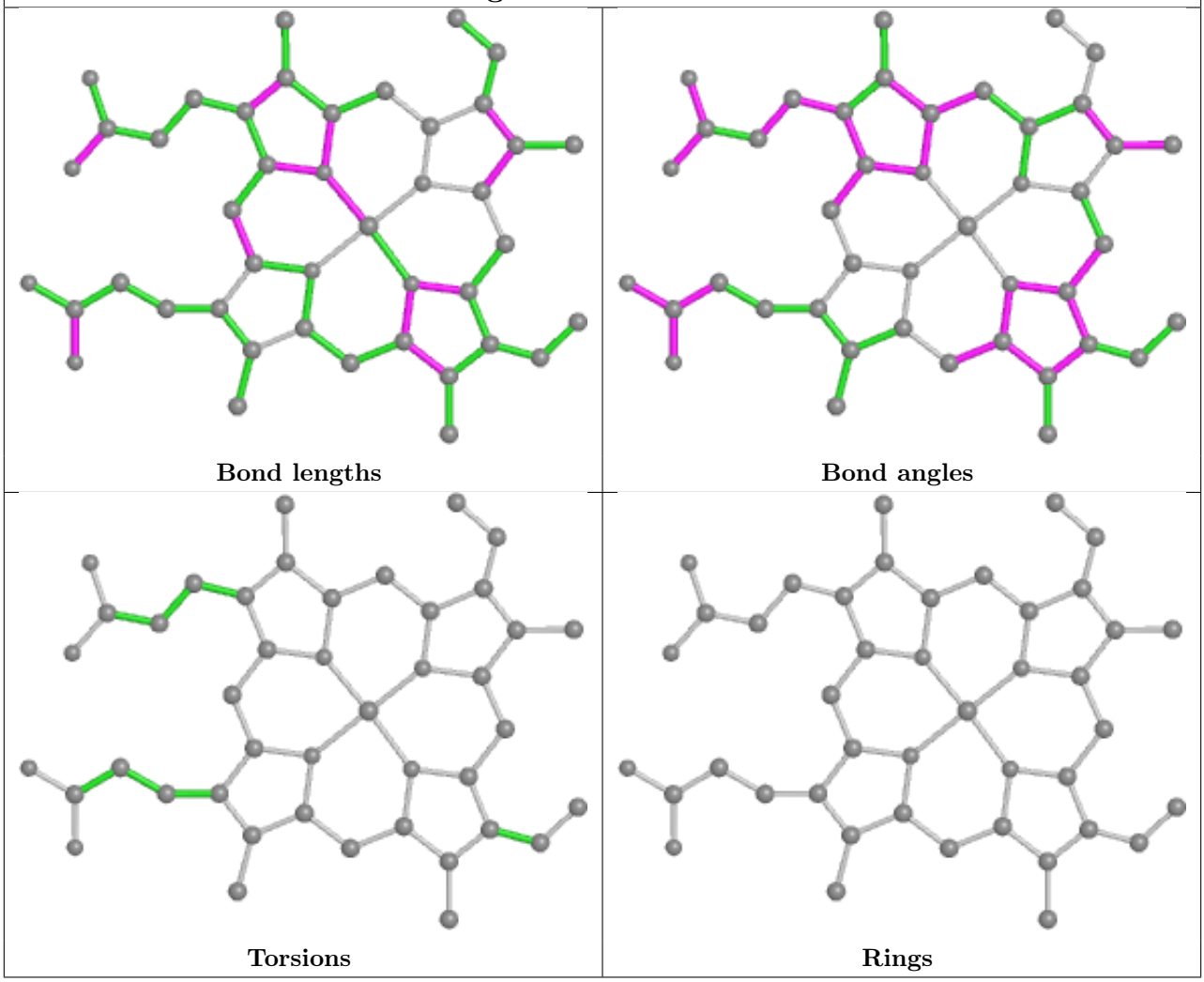
Ligand X2N A 590



Ligand HEM B 580



Ligand HEM A 580



5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

6 Fit of model and data (i)

6.1 Protein, DNA and RNA chains (i)

In the following table, the column labelled '#RSRZ> 2' contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled 'Q< 0.9' lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å^2)	Q<0.9
1	A	484/490 (98%)	0.57	43 (8%) 9 6	36, 63, 130, 189	0
1	B	483/490 (98%)	0.79	61 (12%) 3 2	38, 65, 136, 187	0
All	All	967/980 (98%)	0.68	104 (10%) 5 4	36, 64, 134, 189	0

The worst 5 of 104 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	346	GLY	15.2
1	B	430	ALA	13.4
1	A	346	GLY	12.3
1	B	347	ASP	12.0
1	B	520	TRP	11.0

6.2 Non-standard residues in protein, DNA, RNA chains (i)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates (i)

There are no monosaccharides in this entry.

6.4 Ligands (i)

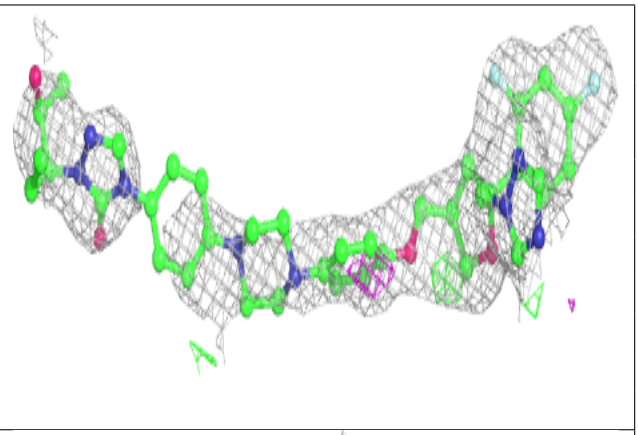
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

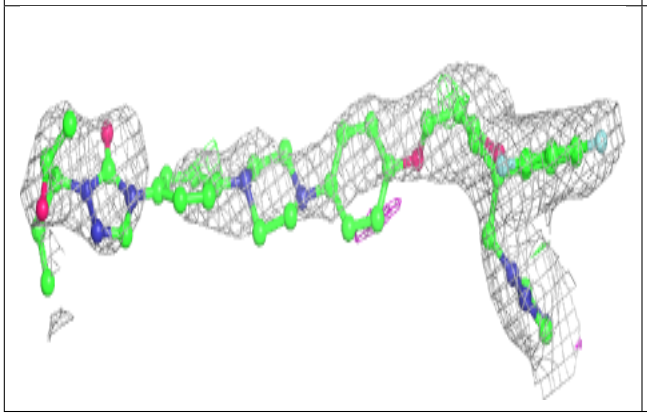
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	X2N	B	590	51/51	0.88	0.32	45,99,118,122	0
3	X2N	A	590	51/51	0.94	0.20	46,74,103,111	0
2	HEM	A	580	43/43	0.97	0.15	34,45,51,53	0
2	HEM	B	580	43/43	0.97	0.17	44,50,55,60	0

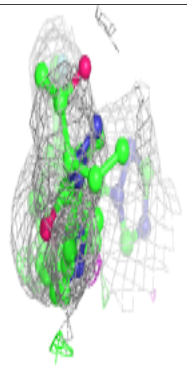
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around X2N B 590:

2mF$_o$-DF$_c$ (at 0.7 rmsd) in gray
mF$_o$-DF$_c$ (at 3 rmsd) in purple (negative) and green (positive)

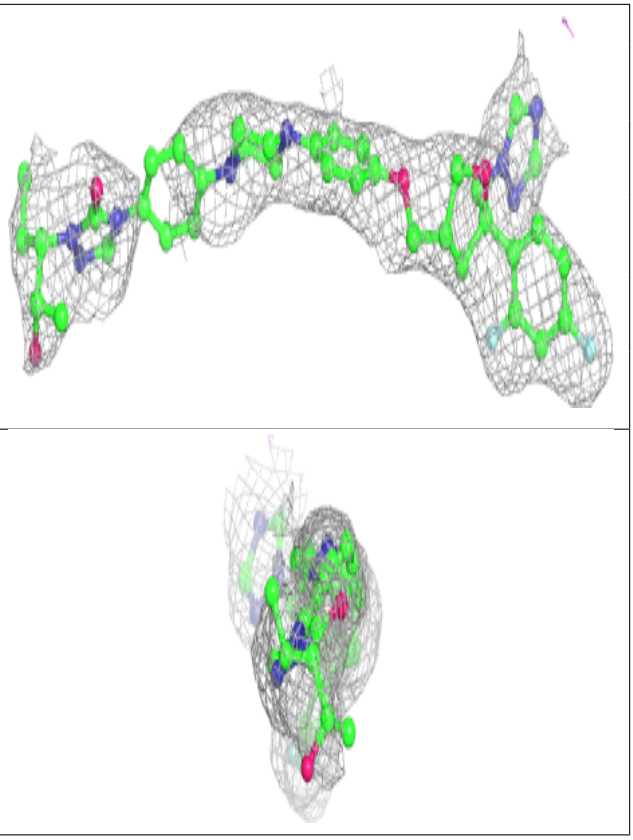






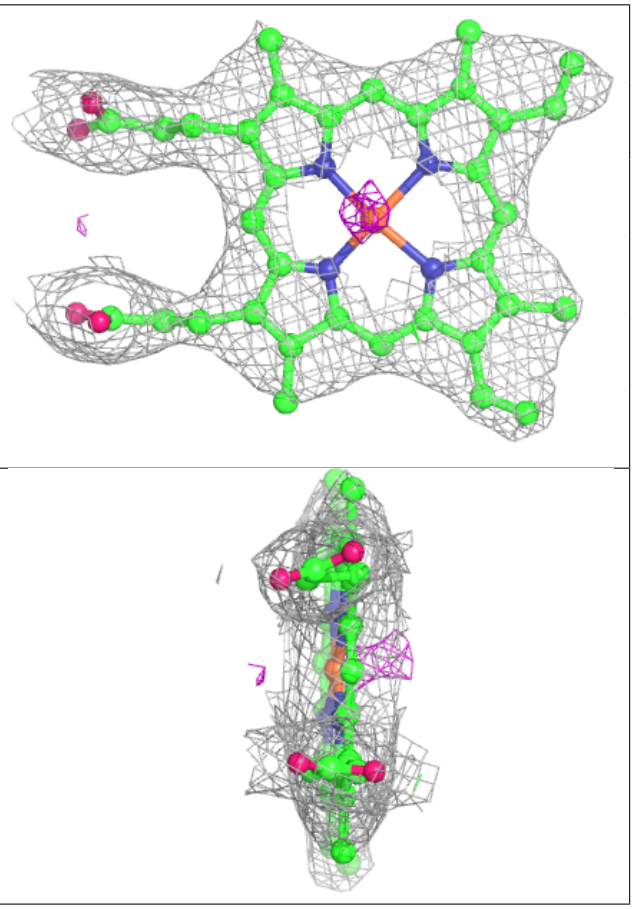
Electron density around X2N A 590:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



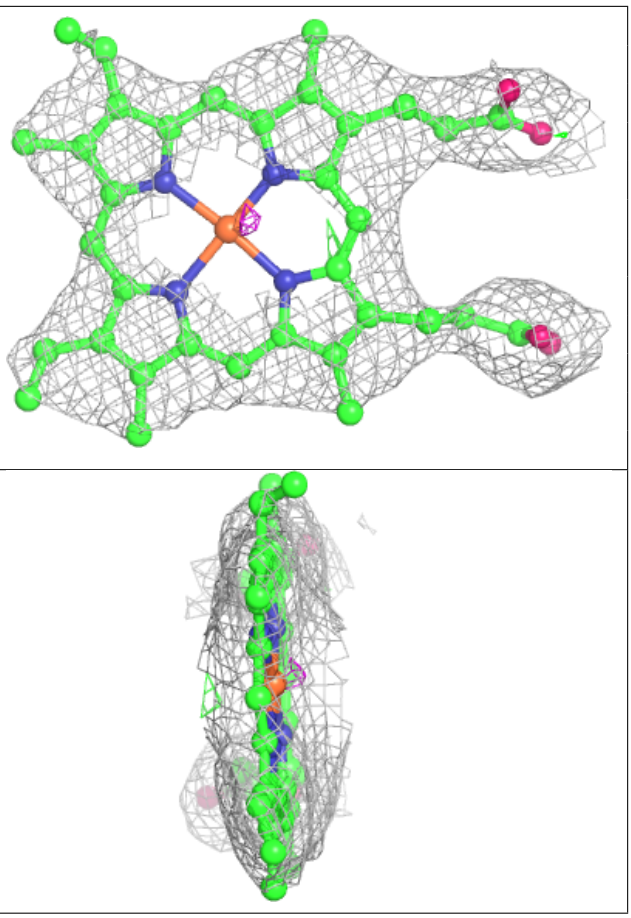
Electron density around HEM A 580:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around HEM B 580:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers (i)

There are no such residues in this entry.